

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN080924\  
 Data File : BN033328.D  
 Acq On : 09 Aug 2024 21:57  
 Operator : MA/JU  
 Sample : P3512-04  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

**Instrument :**  
 BNA\_N  
**ClientSampleId :**  
 927-K1-SD-0.5-1.0-080724

**Manual Integrations**  
**APPROVED**

Reviewed By :Yogesh Patel 08/12/2024  
 Supervised By :mohammad ahmed 08/12/2024

Quant Time: Aug 10 01:10:19 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN080724.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Aug 08 04:39:01 2024  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
<b>Internal Standards</b>							
1) 1,4-Dichlorobenzene-d4	7.580	152	4703	0.400	ng	0.00	
7) Naphthalene-d8	10.349	136	13536	0.400	ng	# 0.00	
13) Acenaphthene-d10	14.212	164	7697	0.400	ng	0.00	
19) Phenanthrene-d10	16.971	188	18787	0.400	ng	# 0.00	
29) Chrysene-d12	21.166	240	17852	0.400	ng	#-0.01	
35) Perylene-d12	23.355	264	20644	0.400	ng	# 0.00	
<b>System Monitoring Compounds</b>							
4) 2-Fluorophenol	5.204	112	2695	0.218	ng	0.00	
5) Phenol-d6	6.757	99	3907	0.231	ng	0.00	
8) Nitrobenzene-d5	8.715	82	2082	0.252	ng	0.00	
11) 2-Methylnaphthalene-d10	11.943	152	4341	0.206	ng	-0.01	
14) 2,4,6-Tribromophenol	15.705	330	633	0.203	ng	-0.01	
15) 2-Fluorobiphenyl	12.844	172	6746	0.233	ng	0.00	
27) Fluoranthene-d10	19.007	212	11448	0.221	ng	0.00	
31) Terphenyl-d14	19.620	244	7958	0.258	ng	-0.01	
<b>Target Compounds</b>							
16) Acenaphthylene	13.934	152	1131	0.030	ng	# 86	Qvalue
17) Acenaphthene	14.277	154	912	0.038	ng	89	
18) Fluorene	15.271	166	1225	0.038	ng	# 95	
25) Phenanthrene	17.008	178	22890	0.421	ng	99	
26) Anthracene	17.095	178	5615	0.106	ng	97	
28) Fluoranthene	19.035	202	74516	1.087	ng	100	
30) Pyrene	19.397	202	63061	0.898	ng	# 95	
32) Benzo(a)anthracene	21.157	228	33167	0.478	ng	98	
33) Chrysene	21.201	228	37725	0.562	ng	97	
34) Bis(2-ethylhexyl)phtha...	21.130	149	1316	0.037	ng	# 92	
36) Indeno(1,2,3-cd)pyrene	25.530	276	24068	0.267	ng	94	
37) Benzo(b)fluoranthene	22.709	252	55948	0.729	ng	97	
38) Benzo(k)fluoranthene	22.747	252	18615m	0.250	ng		
39) Benzo(a)pyrene	23.258	252	34364	0.504	ng	98	
40) Dibenzo(a,h)anthracene	25.542	278	5269	0.074	ng	# 84	
41) Benzo(g,h,i)perylene	26.185	276	25681	0.331	ng	98	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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