

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081724\
 Data File : BN033458.D
 Acq On : 17 Aug 2024 08:27
 Operator : MA/JU
 Sample : P3596-01
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 918-J-WS-081324

Manual Integrations
 APPROVED

Reviewed By :Jagrut
 Upadhyay

Quant Time: Aug 17 23:20:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 16 04:44:11 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							08/19/2024
1) 1,4-Dichlorobenzene-d4	7.559	152	7353	0.400	ng	-0.01	Supervised By :mohammad
7) Naphthalene-d8	10.325	136	20095	0.400	ng	#-0.01	ahmed
13) Acenaphthene-d10	14.199	164	10132	0.400	ng	0.00	
19) Phenanthrene-d10	16.942	188	21230	0.400	ng	-0.01	
29) Chrysene-d12	21.148	240	18526	0.400	ng	0.00	
35) Perylene-d12	23.323	264	19543	0.400	ng	# 0.00	08/21/2024
System Monitoring Compounds							
4) 2-Fluorophenol	5.191	112	2781	0.137	ng	0.00	
5) Phenol-d6	6.743	99	1984	0.072	ng	0.00	
8) Nitrobenzene-d5	8.691	82	5465	0.361	ng	-0.01	
11) 2-Methylnaphthalene-d10	11.918	152	9105	0.294	ng	-0.01	
14) 2,4,6-Tribromophenol	15.688	330	1898	0.365	ng	-0.01	
15) 2-Fluorobiphenyl	12.810	172	14558	0.351	ng	-0.01	
27) Fluoranthene-d10	18.984	212	22664	0.399	ng	0.00	
31) Terphenyl-d14	19.597	244	16877	0.482	ng	0.00	
Target Compounds							Qvalue
18) Fluorene	15.247	166	4595m	0.107	ng		
24) Pentachlorophenol	16.594	266	222	0.039	ng		93
25) Phenanthrene	16.979	178	3065	0.050	ng		99
26) Anthracene	17.078	178	1616	0.030	ng	#	81
28) Fluoranthene	19.012	202	5707	0.076	ng		99
30) Pyrene	19.379	202	3893	0.053	ng	#	94
32) Benzo(a)anthracene	21.139	228	1866	0.027	ng	#	83
33) Chrysene	21.184	228	2774	0.040	ng	#	91
34) Bis(2-ethylhexyl)phtha...	21.104	149	4003	0.127	ng	#	96
36) Indeno(1,2,3-cd)pyrene	25.478	276	1904	0.024	ng		99
37) Benzo(b)fluoranthene	22.677	252	2998	0.041	ng	#	18
38) Benzo(k)fluoranthene	22.721	252	1763	0.023	ng	#	1
39) Benzo(a)pyrene	23.227	252	1318	0.021	ng	#	1
41) Benzo(g,h,i)perylene	26.139	276	1763	0.025	ng	#	55

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

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