

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120924\
 Data File : BN035502.D
 Acq On : 09 Dec 2024 12:16
 Operator : RC/JU
 Sample : P5142-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TOWER-1

Manual Integrations
 APPROVED

Reviewed By :Yogesh
 Patel

Quant Time: Dec 09 13:27:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN112724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 27 23:03:24 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							12/11/2024
1) 1,4-Dichlorobenzene-d4	7.300	152	2170	0.400	ng	0.00	Supervised By :mohammad Ahmed
7) Naphthalene-d8	10.041	136	5734	0.400	ng	-0.01	
13) Acenaphthene-d10	13.956	164	4922	0.400	ng	-0.01	
19) Phenanthrene-d10	16.723	188	11976	0.400	ng	#-0.01	
29) Chrysene-d12	20.964	240	11160	0.400	ng	# 0.00	
35) Perylene-d12	23.055	264	10129	0.400	ng	-0.01	12/11/2024
System Monitoring Compounds							
4) 2-Fluorophenol	4.960	112	704	0.130	ng	0.00	
5) Phenol-d6	6.506	99	579	0.089	ng	0.00	
8) Nitrobenzene-d5	8.429	82	1666m	0.476	ng	-0.01	
11) 2-Methylnaphthalene-d10	11.641	152	3824	0.426	ng	-0.02	
14) 2,4,6-Tribromophenol	15.464	330	1258	0.360	ng	-0.01	
15) 2-Fluorobiphenyl	12.564	172	6512	0.350	ng	-0.01	
27) Fluoranthene-d10	18.775	212	15399	0.454	ng	0.00	
31) Terphenyl-d14	19.402	244	12062	0.548	ng	0.00	
Target Compounds							Qvalue
2) 1,4-Dioxane	2.996	88	410	0.198	ng	#	22
9) Naphthalene	10.094	128	520	0.034	ng	#	64
24) Pentachlorophenol	16.375	266	356	0.099	ng	#	88
25) Phenanthrene	16.760	178	776	0.024	ng	#	66
28) Fluoranthene	18.808	202	1072	0.024	ng	#	89
30) Pyrene	19.175	202	1041	0.025	ng	#	94
32) Benzo(a)anthracene	20.946	228	1331	0.034	ng	#	59
33) Chrysene	21.000	228	1514	0.038	ng	#	66
34) Bis(2-ethylhexyl)phtha...	20.929	149	6307	0.409	ng	#	68
36) Indeno(1,2,3-cd)pyrene	25.090	276	1211	0.031	ng	#	98
37) Benzo(b)fluoranthene	22.444	252	1616	0.044	ng	#	16
38) Benzo(k)fluoranthene	22.485	252	1332	0.037	ng	#	4
39) Benzo(a)pyrene	22.965	252	825	0.027	ng	#	1
40) Dibenzo(a,h)anthracene	25.105	278	825	0.026	ng	#	12
41) Benzo(g,h,i)perylene	25.707	276	968	0.030	ng	#	56

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

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