

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040524\
 Data File : BN030666.D
 Acq On : 04 Apr 2024 18:17
 Operator : MA/JU
 Sample : P1964-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RSB-02-47-040224

Quant Time: Apr 04 18:46:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN033024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 30 00:04:36 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.697	152	7489	0.400	ng	0.00
7) Naphthalene-d8	10.475	136	22046	0.400	ng	-0.01
13) Acenaphthene-d10	14.327	164	13509	0.400	ng	-0.01
19) Phenanthrene-d10	17.078	188	30840	0.400	ng	#-0.01
29) Chrysene-d12	21.276	240	26643	0.400	ng	# 0.00
35) Perylene-d12	23.521	264	23766	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.284	112	2462	0.151	ng	-0.01
5) Phenol-d6	6.866	99	1870	0.089	ng	0.00
8) Nitrobenzene-d5	8.841	82	4469	0.288	ng	-0.01
11) 2-Methylnaphthalene-d10	12.070	152	8917	0.264	ng	-0.01
14) 2,4,6-Tribromophenol	15.824	330	1877	0.368	ng	-0.01
15) 2-Fluorobiphenyl	12.948	172	11819	0.226	ng	-0.02
27) Fluoranthene-d10	19.117	212	25383	0.305	ng	0.00
31) Terphenyl-d14	19.721	244	18529	0.301	ng	0.00
Target Compounds						Qvalue
24) Pentachlorophenol	16.730	266	30	0.175	ng	# 47
32) Benzo(a)anthracene	21.267	228	2538	0.027	ng	# 90
33) Chrysene	21.312	228	2579	0.025	ng	# 91
34) Bis(2-ethylhexyl)phtha...	21.205	149	5571	0.136	ng	# 98
37) Benzo(b)fluoranthene	22.846	252	2312	0.024	ng	# 12
38) Benzo(k)fluoranthene	22.890	252	2298	0.024	ng	# 13

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

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