

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110423\
 Data File : BN028477.D
 Acq On : 04 Nov 2023 16:35
 Operator : MA/JU
 Sample : 05242-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SP-303-20231102

Quant Time: Nov 06 04:46:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN110123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 02 06:26:50 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.717	152	9835	0.400	ng	0.00
7) Naphthalene-d8	10.509	136	30041	0.400	ng	# 0.00
13) Acenaphthene-d10	14.363	164	18587	0.400	ng	0.00
19) Phenanthrene-d10	17.109	188	42208	0.400	ng	#-0.01
29) Chrysene-d12	21.328	240	33173	0.400	ng	0.00
35) Perylene-d12	23.646	264	30127	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.298	112	153	0.006	ng	0.00
5) Phenol-d6	6.887	99	158	0.005	ng	0.00
8) Nitrobenzene-d5	8.865	82	6193	0.319	ng	-0.01
11) 2-Methylnaphthalene-d10	12.106	152	13973	0.325	ng	0.00
14) 2,4,6-Tribromophenol	15.855	330	403	0.056	ng	0.00
15) 2-Fluorobiphenyl	13.006	172	2665	0.121	ng	0.00
27) Fluoranthene-d10	19.157	212	40307	0.386	ng	0.00
31) Terphenyl-d14	19.765	244	29228	0.418	ng	0.00
Target Compounds						
34) Bis(2-ethylhexyl)phtha...	21.266	149	17908	0.263	ng	# 99
36) Indeno(1,2,3-cd)pyrene	26.026	276	1146	0.057	ng	# 79
37) Benzo(b)fluoranthene	22.944	252	1434	0.079	ng	# 1

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

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