

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110423\
 Data File : BN028478.D
 Acq On : 04 Nov 2023 17:12
 Operator : MA/JU
 Sample : 05242-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SP-201-20231102

Quant Time: Nov 06 04:46:57 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	9654	0.400	ng	-0.01
7) Naphthalene-d8	10.508	136	29411	0.400	ng	# 0.00
13) Acenaphthene-d10	14.363	164	18217	0.400	ng	0.00
19) Phenanthrene-d10	17.108	188	42635	0.400	ng	#-0.01
29) Chrysene-d12	21.328	240	34455	0.400	ng	0.00
35) Perylene-d12	23.645	264	32664	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.297	112	2681	0.106	ng	0.00
5) Phenol-d6	6.886	99	1900	0.059	ng	0.00
8) Nitrobenzene-d5	8.864	82	5702	0.300	ng	-0.01
11) 2-Methylnaphthalene-d10	12.102	152	12611	0.300	ng	-0.01
14) 2,4,6-Tribromophenol	15.854	330	2587	0.364	ng	0.00
15) 2-Fluorobiphenyl	13.005	172	2279	0.105	ng	0.00
27) Fluoranthene-d10	19.156	212	39022	0.370	ng	0.00
31) Terphenyl-d14	19.765	244	30686	0.423	ng	0.00
Target Compounds						
34) Bis(2-ethylhexyl)phtha...	21.265	149	5884	0.083	ng	# 99
36) Indeno(1,2,3-cd)pyrene	26.025	276	1043	0.056	ng	# 74
37) Benzo(b)fluoranthene	22.946	252	1182	0.077	ng	# 1

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

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