

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101123\
 Data File : BN028240.D
 Acq On : 11 Oct 2023 19:38
 Operator : MA/JU
 Sample : 04819-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SW-AP-01-100923

Quant Time: Oct 12 04:05:54 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 12 04:03:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	3109	0.400	ng	0.00
7) Naphthalene-d8	10.735	136	10739	0.400	ng	# 0.00
13) Acenaphthene-d10	14.566	164	7049	0.400	ng	0.00
19) Phenanthrene-d10	17.319	188	15511	0.400	ng	# 0.00
29) Chrysene-d12	21.524	240	14521	0.400	ng	0.00
35) Perylene-d12	23.996	264	15072	0.400	ng	0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.459	112	1203	0.130	ng	0.00
5) Phenol-d6	7.106	99	852	0.073	ng	0.01
8) Nitrobenzene-d5	9.123	82	2439	0.267	ng	0.00
11) 2-Methylnaphthalene-d10	12.324	152	4731	0.297	ng	0.00
14) 2,4,6-Tribromophenol	16.065	330	1185	0.367	ng	-0.01
15) 2-Fluorobiphenyl	13.187	172	7159	0.282	ng	0.00
27) Fluoranthene-d10	19.356	212	14888	0.382	ng	0.00
31) Terphenyl-d14	19.946	244	12286	0.363	ng	0.00
Target Compounds						
28) Fluoranthene	19.388	202	1476	0.031	ng	Qvalue 98
30) Pyrene	19.755	202	1300	0.024	ng	93
32) Benzo(a)anthracene	21.506	228	1556	0.031	ng	# 62
33) Chrysene	21.560	228	2034	0.043	ng	# 67
34) Bis(2-ethylhexyl)phtha...	21.372	149	10254	0.316	ng	100
36) Indeno(1,2,3-cd)pyrene	26.595	276	1535	0.030	ng	# 91
37) Benzo(b)fluoranthene	23.230	252	2846	0.058	ng	# 1
41) Benzo(g,h,i)perylene	27.411	276	1545	0.036	ng	# 25

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

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