

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032122\
 Data File : BP009473.D
 Acq On : 21 Mar 2022 16:38
 Operator : CG/JU
 Sample : N1983-01DL 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 124035DL

Quant Time: Mar 21 17:16:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 17 18:26:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.792	152	414734	20.000	ng	0.00
21) Naphthalene-d8	10.586	136	1633905	20.000	ng	0.00
39) Acenaphthene-d10	14.433	164	1062203	20.000	ng	-0.01
64) Phenanthrene-d10	17.198	188	2272653	20.000	ng	0.00
76) Chrysene-d12	21.309	240	2423506	20.000	ng	0.00
86) Perylene-d12	23.715	264	2581608	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.398	112	502385	21.149	ng	0.00
7) Phenol-d6	6.987	99	615963	19.174	ng	0.00
23) Nitrobenzene-d5	8.951	82	379848	12.354	ng	0.00
42) 2,4,6-Tribromophenol	15.933	330	263063	15.008	ng	-0.01
45) 2-Fluorobiphenyl	13.057	172	888015	11.590	ng	0.00
79) Terphenyl-d14	19.774	244	1603533	11.876	ng	0.00
Target Compounds						
						Qvalue
37) 2-Methylnaphthalene	12.257	142	266031	4.505	ng	98
38) 1-Methylnaphthalene	12.475	142	193206	3.359	ng	99
52) Acenaphthene	14.498	154	656680	11.414	ng	100
55) Dibenzofuran	14.839	168	793152	8.206	ng	99
58) Fluorene	15.492	166	886665	11.665	ng	98
71) Phenanthrene	17.239	178	3958632	32.540	ng	100
72) Anthracene	17.333	178	501957	4.179	ng	98
75) Fluoranthene	19.221	202	2174105	14.761	ng	99
78) Pyrene	19.574	202	1436410	8.995	ng	100
81) Benzo(a)anthracene	21.298	228	423522	2.660	ng	99
83) Chrysene	21.345	228	381353	2.529	ng	98

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

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