

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020824\
 Data File : BP019621.D
 Acq On : 08 Feb 2024 22:16
 Operator : MA/JU
 Sample : P1413-01 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VNJ-209

Quant Time: Feb 09 00:44:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 00:51:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	101700	20.000	ng	0.00
21) Naphthalene-d8	10.710	136	387699	20.000	ng	0.00
39) Acenaphthene-d10	14.540	164	245310	20.000	ng	-0.01
64) Phenanthrene-d10	17.298	188	507634	20.000	ng	0.00
76) Chrysene-d12	21.392	240	450567	20.000	ng	0.00
86) Perylene-d12	23.816	264	517960	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	338502	53.652	ng	0.00
7) Phenol-d6	7.081	99	472473	55.538	ng	0.00
23) Nitrobenzene-d5	9.075	82	304468	34.029	ng	-0.01
42) 2,4,6-Tribromophenol	16.039	330	225464	62.949	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	735548	37.068	ng	-0.01
79) Terphenyl-d14	19.863	244	1100773	40.178	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.339	178	62828	2.352	ng	96
75) Fluoranthene	19.310	202	107764	3.302	ng	98
78) Pyrene	19.663	202	84064	2.644	ng	97

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

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