

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
 Data File : BP020084.D
 Acq On : 22 Apr 2024 20:22
 Operator : MA/JU
 Sample : P2104-21
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COAL8

Quant Time: Apr 23 01:36:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 23 01:17:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.640	152	144655	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	567390	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164	297936	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.163	188	517691	20.000	ng/ul	0.00
79) Chrysene-d12	21.663	240	484697	20.000	ng/ul	0.00
88) Perylene-d12	25.062	264	559391	20.000	ng/ul	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	11673	3.306	ng/uL	0.00
4) Pyridine-d5	3.646	84	26238	2.525	ng/ul	0.01
7) Phenol-d5	6.863	99	1142	0.091	ng/ul	0.04
9) Bis-(2-Chloroethyl)eth...	6.993	67	188411	21.381	ng/ul	0.00
11) 2-Chlorophenol-d4	7.187	132	3465	0.403	ng/ul	0.01
15) 4-Methylphenol-d8	8.357	113	4988	0.506	ng/ul	0.01
21) Nitrobenzene-d5	8.799	128	97920	23.992	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	101681	23.048	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	675	0.077	ng/ul	0.06
31) 4-Chloroaniline-d4	10.610	131	417	0.036	ng/ul	0.04
46) Dimethylphthalate-d6	13.734	166	560123	27.509	ng/ul	0.00
49) Acenaphthylene-d8	13.998	160	3764	0.155	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143	601	0.194	ng/ul	-0.01
60) Fluorene-d10	15.345	176	49345	2.722	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.504	200	13955	4.614	ng/ul	0.02
73) Anthracene-d10	17.163	188	517691	22.321	ng/ul	-0.10
81) Pyrene-d10	19.663	212	5396	0.207	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.827	264	3114	0.114	ng/ul	0.02

Target Compounds	Qvalue

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

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