

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
 Data File : BP013744.D
 Acq On : 03 Feb 2023 17:14
 Operator : CG/JU
 Sample : 01381-02
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A44M6

Quant Time: Feb 03 23:52:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 03 23:44:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.134	152	217018	20.000	ng/ul	0.00
20) Naphthalene-d8	10.957	136	922749	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.769	164	670326	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.522	188	1685176	20.000	ng/ul	0.00
79) Chrysene-d12	21.598	240	1944600	20.000	ng/ul	0.00
88) Perylene-d12	24.169	264	2293275	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.475	96	27124	5.299	ng/uL	0.00
4) Pyridine-d5	3.917	84	129173	8.795	ng/ul	0.00
7) Phenol-d5	7.275	99	99719	5.228	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.446	67	329736	29.476	ng/ul	0.00
11) 2-Chlorophenol-d4	7.658	132	288362	20.624	ng/ul	0.00
15) 4-Methylphenol-d8	8.834	113	191794	12.614	ng/ul	0.00
21) Nitrobenzene-d5	9.305	128	179510	30.495	ng/ul	0.00
24) 2-Nitrophenol-d4	10.034	143	191495	27.914	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.569	165	398364	24.832	ng/ul	0.00
31) 4-Chloroaniline-d4	11.093	131	461963	21.804	ng/ul	0.00
46) Dimethylphthalate-d6	14.169	166	1711340	32.388	ng/ul	0.00
49) Acenaphthylene-d8	14.463	160	1689681	28.943	ng/ul	0.00
54) 4-Nitrophenol-d4	14.957	143	24163	2.713	ng/ul	0.00
60) Fluorene-d10	15.757	176	1511531	33.188	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.869	200	140649	14.894	ng/ul	0.00
73) Anthracene-d10	17.622	188	2686297	34.198	ng/ul	0.00
81) Pyrene-d10	19.839	212	3537384	32.615	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.004	264	4039041	34.501	ng/ul	0.00

Target Compounds Qvalue

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

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