

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110723\
 Data File : BP017970.D
 Acq On : 08 Nov 2023 19:41
 Operator : MA/JU
 Sample : 05133-01
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A49Z5

Quant Time: Nov 08 21:44:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 07 03:40:56 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	74354	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.669	136	308586	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.516	164	199189	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	449342	20.000	ng/ul	0.00
79) Chrysene-d12	21.410	240	423271	20.000	ng/ul	0.00
88) Perylene-d12	23.892	264	478116	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	6434	3.312	ng/uL	0.00
4) Pyridine-d5	3.669	84	157	0.029	ng/ul	0.00
7) Phenol-d5	7.028	99	20805	2.984	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.204	67	46632	11.087	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	48366	9.027	ng/ul	0.00
15) 4-Methylphenol-d8	8.581	113	31098	5.396	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	28284	11.469	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	30170	10.622	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	46497	8.930	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	25673	3.402	ng/ul	0.01
46) Dimethylphthalate-d6	13.939	166	225011	12.883	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	219593	11.458	ng/ul	0.00
54) 4-Nitrophenol-d4	14.833	143	1781	0.663	ng/ul	0.06
60) Fluorene-d10	15.516	176	181625	12.424	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	26603	8.096	ng/ul	0.00
73) Anthracene-d10	17.392	188	296973	12.711	ng/ul	0.00
81) Pyrene-d10	19.639	212	375541	13.977	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	356193	13.333	ng/ul	0.00

Target Compounds Qvalue

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

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