

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110723\
 Data File : BP017974.D
 Acq On : 08 Nov 2023 22:31
 Operator : MA/JU
 Sample : PB156654BL
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK654

Quant Time: Nov 08 23:24:09 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	7.858	152	51805	20.000	ng/u1	0.00
20) Naphthalene-d8	10.669	136	215052	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.516	164	135497	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.287	188	314550	20.000	ng/u1	-0.01
79) Chrysene-d12	21.410	240	283906	20.000	ng/u1	0.00
88) Perylene-d12	23.886	264	307734	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	11810	8.725	ng/uL	0.00
4) Pyridine-d5	3.675	84	143121	38.099	ng/u1	0.00
7) Phenol-d5	7.022	99	178567	36.757	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	113982	38.895	ng/u1	0.00
11) 2-Chlorophenol-d4	7.375	132	147228	39.440	ng/u1	-0.01
15) 4-Methylphenol-d8	8.575	113	142929	35.596	ng/u1	0.00
21) Nitrobenzene-d5	9.058	128	71694	41.717	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	81013	40.927	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	134553	37.079	ng/u1	0.00
31) 4-Chloroaniline-d4	10.846	131	190916	36.306	ng/u1	0.00
46) Dimethylphthalate-d6	13.940	166	498098	41.924	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	514924	39.497	ng/u1	0.00
54) 4-Nitrophenol-d4	14.775	143	56121	30.722	ng/u1	0.00
60) Fluorene-d10	15.516	176	406004	40.828	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	70113	30.480	ng/u1	-0.01
73) Anthracene-d10	17.387	188	668142	40.853	ng/u1	-0.01
81) Pyrene-d10	19.639	212	791590	43.923	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.721	264	744829	43.317	ng/u1	-0.01

Target Compounds Qvalue

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

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