

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091324\
 Data File : BP021907.D
 Acq On : 13 Sep 2024 13:47
 Operator : RC/JU
 Sample : PB163354BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK354

Quant Time: Sep 14 04:53:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 12 12:48:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	137003	20.000	ng/u1	0.03
20) Naphthalene-d8	10.528	136	546446	20.000	ng/u1	0.02
38) Acenaphthene-d10	14.369	164	399645	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.157	188	927166	20.000	ng/u1	0.00
79) Chrysene-d12	21.586	240	955441	20.000	ng/u1	0.00
88) Perylene-d12	24.915	264	1035448	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	19579	4.623	ng/uL	0.01
4) Pyridine-d5	3.681	84	221767	17.784	ng/u1	0.02
7) Phenol-d5	6.934	99	306254	26.161	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.099	67	195092	28.693	ng/u1	0.03
11) 2-Chlorophenol-d4	7.299	132	244099	29.955	ng/u1	0.02
15) 4-Methylphenol-d8	8.457	113	182665	19.544	ng/u1	0.03
21) Nitrobenzene-d5	8.899	128	91178	26.791	ng/u1	0.02
24) 2-Nitrophenol-d4	9.616	143	97496	25.200	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.151	165	255024	29.766	ng/u1	0.02
31) 4-Chloroaniline-d4	10.657	131	336157	26.923	ng/u1	0.02
46) Dimethylphthalate-d6	13.769	166	918992	30.608	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	938887	28.507	ng/u1	0.00
54) 4-Nitrophenol-d4	14.569	143	136219	24.000	ng/u1	0.00
60) Fluorene-d10	15.369	176	785280	27.516	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	119428	22.515	ng/u1	0.00
73) Anthracene-d10	17.257	188	1297911	30.136	ng/u1	0.00
81) Pyrene-d10	19.621	212	1633749	30.906	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.692	264	1632728	30.299	ng/u1	0.00

Target Compounds	Qvalue

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

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