

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011960.D
 Acq On : 06 Oct 2022 09:20
 Operator : CG/JU
 Sample : PB148128BL
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 06 23:04:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 05 23:18:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	265177	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	1111454	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	672244	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1408907	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1318422	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1334429	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	44187	7.349	ng/uL	0.00
4) Pyridine-d5	3.717	84	547288	31.075	ng/ul	0.00
7) Phenol-d5	7.040	99	749131	32.724	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	428581	33.775	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	618895	34.247	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	584280	30.875	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	295140	34.863	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	308264	34.034	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	556782	33.234	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	824204	32.270	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	1661919	34.571	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1931139	35.238	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	290880	29.220	ng/ul	0.00
60) Fluorene-d10	15.522	176	1503435	36.012	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	223192	26.173	ng/ul	0.00
73) Anthracene-d10	17.381	188	2414570	37.623	ng/ul	0.00
81) Pyrene-d10	19.616	212	2749455	41.178	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2673415	40.021	ng/ul	0.00

Target Compounds Qvalue

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

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