

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011885.D
 Acq On : 03 Oct 2022 23:04
 Operator : CG/JU
 Sample : PB148012BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK012

Quant Time: Oct 04 00:50:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 01 00:53:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	256379	20.000	ng/u1	0.00
20) Naphthalene-d8	10.699	136	1072324	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.534	164	653715	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.286	188	1396081	20.000	ng/u1	-0.01
79) Chrysene-d12	21.374	240	1396188	20.000	ng/u1	-0.01
88) Perylene-d12	23.810	264	1424366	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	39920	6.867	ng/uL	0.00
4) Pyridine-d5	3.728	84	396516	23.287	ng/u1	0.00
7) Phenol-d5	7.052	99	694503	31.378	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	393453	32.071	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.416	132	592000	33.883	ng/u1	0.00
15) 4-Methylphenol-d8	8.599	113	559880	30.600	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	280362	34.325	ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	302770	34.647	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	535423	33.126	ng/u1	0.00
31) 4-Chloroaniline-d4	10.840	131	459497	18.647	ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	1614228	34.531	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	1930222	36.220	ng/u1	0.00
54) 4-Nitrophenol-d4	14.734	143	298782	30.865	ng/u1	0.00
60) Fluorene-d10	15.528	176	1444331	35.577	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	231675	27.417	ng/u1	0.00
73) Anthracene-d10	17.386	188	2240910	35.238	ng/u1	-0.01
81) Pyrene-d10	19.622	212	2712667	38.364	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.651	264	2699403	37.859	ng/u1	-0.01

Target Compounds	Qvalue

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

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