

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012239.D
 Acq On : 21 Oct 2022 11:29
 Operator : CG/JU
 Sample : PB148371BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK371

Quant Time: Oct 21 22:30:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 00:51:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	401711	20.000	ng/u1	0.00
20) Naphthalene-d8	10.628	136	1723314	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.475	164	1023843	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.234	188	1844443	20.000	ng/u1	0.00
79) Chrysene-d12	21.322	240	1431057	20.000	ng/u1	-0.01
88) Perylene-d12	23.727	264	1506943	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	58516	5.895	ng/uL	0.00
4) Pyridine-d5	3.676	84	813224	28.432	ng/u1	0.00
7) Phenol-d5	6.993	99	1056112	30.453	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	640596	30.458	ng/u1	0.00
11) 2-Chlorophenol-d4	7.352	132	833140	31.561	ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	830602	30.816	ng/u1	-0.01
21) Nitrobenzene-d5	8.993	128	415835	32.599	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	450580	32.739	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.252	165	781683	30.280	ng/u1	0.00
31) 4-Chloroaniline-d4	10.775	131	1144469	31.203	ng/u1	0.00
46) Dimethylphthalate-d6	13.887	166	2215434	31.107	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	2554714	31.426	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.687	143	276344	20.711	ng/u1	0.00
60) Fluorene-d10	15.469	176	1881184	30.703	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	138023	13.154	ng/u1	0.00
73) Anthracene-d10	17.334	188	2610403	32.288	ng/u1	0.00
81) Pyrene-d10	19.569	212	2577813	31.638	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.574	264	2499563	33.076	ng/u1	-0.01

Target Compounds Qvalue

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

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