

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012430.D
 Acq On : 01 Nov 2022 16:43
 Operator : CG/JU
 Sample : PB148596BL
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK596

Quant Time: Nov 01 23:00:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 01 01:03:58 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	324448	20.000	ng/u1	0.00
20) Naphthalene-d8	10.593	136	1483190	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.439	164	985684	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.204	188	2156710	20.000	ng/u1	0.00
79) Chrysene-d12	21.298	240	2155972	20.000	ng/u1	0.00
88) Perylene-d12	23.686	264	1914595	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	46530	5.775	ng/uL	0.00
4) Pyridine-d5	3.658	84	601155	25.971	ng/u1	0.00
7) Phenol-d5	6.963	99	836023	28.729	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	513259	29.545	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	654012	30.776	ng/u1	0.00
15) 4-Methylphenol-d8	8.510	113	676099	29.375	ng/u1	0.00
21) Nitrobenzene-d5	8.957	128	349810	36.370	ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	381725	38.237	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	639819	29.573	ng/u1	0.00
31) 4-Chloroaniline-d4	10.740	131	959171	29.775	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	2200549	32.613	ng/u1	0.00
49) Acenaphthylene-d8	14.134	160	2467123	31.995	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	384239	30.088	ng/u1	0.00
60) Fluorene-d10	15.439	176	1914171	33.134	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	321190	28.604	ng/u1	0.00
73) Anthracene-d10	17.304	188	3069658	33.033	ng/u1	0.00
81) Pyrene-d10	19.545	212	3600268	33.261	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	3099380	33.131	ng/u1	-0.01

Target Compounds	Qvalue

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

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