

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
 Data File : BP012743.D
 Acq On : 26 Nov 2022 00:32
 Operator : CG/JU
 Sample : N5592-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0AA2

Quant Time: Nov 26 07:18:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 25 15:59:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.022	152	571650	20.000	ng/u1	0.00
20) Naphthalene-d8	10.840	136	2538883	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.657	164	1665721	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.410	188	3410380	20.000	ng/u1	0.00
79) Chrysene-d12	21.498	240	2243878	20.000	ng/u1	0.00
88) Perylene-d12	24.015	264	1801348	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	61681	4.624	ng/uL	0.00
4) Pyridine-d5	3.822	84	349222	8.748	ng/u1	0.00
7) Phenol-d5	7.169	99	303557	6.462	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.346	67	846195	29.234	ng/u1	0.00
11) 2-Chlorophenol-d4	7.546	132	817487	22.839	ng/u1	0.00
15) 4-Methylphenol-d8	8.722	113	583272	15.411	ng/u1	0.00
21) Nitrobenzene-d5	9.187	128	496632	31.960	ng/u1	0.00
24) 2-Nitrophenol-d4	9.916	143	506570	30.701	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	1025073	26.626	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	958613	18.183	ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	3869925	33.753	ng/u1	0.00
49) Acenaphthylene-d8	14.351	160	4236477	31.668	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	77295	3.781	ng/u1	-0.01
60) Fluorene-d10	15.651	176	3338523	33.249	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	406675	24.950	ng/u1	0.00
73) Anthracene-d10	17.510	188	5346060	35.612	ng/u1	0.00
81) Pyrene-d10	19.739	212	5454177	41.832	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.851	264	3216607	36.354	ng/u1	0.00

Target Compounds Qvalue

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

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