

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013064.D
 Acq On : 11 Dec 2022 05:31
 Operator : CG/JU
 Sample : N5927-07
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 12 03:31:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 12 01:32:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	611319	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	2674791	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.610	164	1814662	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	4188337	20.000	ng/ul	0.00
79) Chrysene-d12	21.445	240	3893122	20.000	ng/ul	0.00
88) Perylene-d12	23.939	264	3909182	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	66422	4.646	ng/uL	0.00
4) Pyridine-d5	3.781	84	375608	9.248	ng/ul	0.00
7) Phenol-d5	7.122	99	335791	6.744	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	1047440	34.871	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132	1000482	25.524	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	656776	16.098	ng/ul	0.00
21) Nitrobenzene-d5	9.134	128	700973	35.669	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	737484	33.333	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	1250446	29.097	ng/ul	0.00
31) 4-Chloroaniline-d4	10.916	131	1704963	28.103	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	5123032	36.733	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	5529966	36.085	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	122670	4.890	ng/ul	0.00
60) Fluorene-d10	15.604	176	4424361	37.498	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	800515	29.701	ng/ul	0.00
73) Anthracene-d10	17.463	188	7288091	38.570	ng/ul	0.00
81) Pyrene-d10	19.692	212	8682314	38.853	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.774	264	7855200	40.314	ng/ul	0.00

Target Compounds Qvalue

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

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